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Catalog Handbook of Fine Chemicals 1996-1997





United States



-	ll IIIIIIIIIII		
■ Acety	i'chi ■		\$
_	Acetyl chloride, 98.5 + %, A.C.S. reagent [75-36-5] CH₃COCI	25mL 500mL x500mL 2.5L†	13.20 27.00 129.60 106.95
11,418-9	Evapn. residue ≤0.005% Fe ≤5 ppm Insolubles ≤0.0025% Acetyl chloride, 98% [75-36-5] CH₃COCI	25g 500g 1kg	11.50 20.15 29.20
32.012-9	Acetyl chloride, 98% [75-36-5] CH ₃ COCI	_	44.55
24,707-3	(Packaged in poly-coated bottle) 2-Acetyl-5-chlorothlophene, 99% [6310-09-4] (5-chloro-2-thienyl methylketone) FW 160.62 mp 46-49° bp 117-118°/17mm Fp.227°F(108°C) Beil. 17,287 FT-NMR 1(3),53C FT-IR 1(2),597D SI 368,A,2 R&S 1(2),2337B RTECS# OB1745000	. 5g 25g	20.30 67.70
85,968-0 *	Acetylcholine bromide, 98% [66-23-9] CH ₂ CO ₂ CH ₂ CH ₃ N(CH ₃),Br FW 226.12	25g 100g	13.30 41.90
*	Acetylcholine chloride, 98% [60-31-1] CH,CO,CH,CH,N(CH,),CI FW 181.66	500g	11.10 38.30 172.35 20.30
*	mp 162-164° FT-IR 1(1),6780 ST 120,D, 'asiety 2,508 Title (1),6780 ST 120,D, 'asiety 2,508 Title		21.30 70.70
21,407·1	Acetyl cyanide, see 26,921-2, Pyruvonitrile page 1289 1-Acetyl-2-(cyanoacetyl)hydrazine, see 38,024-5, Cyanoacetic 2-acetylhydrazide page 408 N.Acetyl-2-cyanoalycine ethyl ester, see E960-9, Ethyl	25g	70.70
15,649-3	acetamidocyanoacetate page 667 2-Acetyl-1,3-cyclohexanedione, 98% [4056-73-9] CH ₃ COC ₆ H ₇ (= O), FW 154.17 mp 20° bp 85 %0.1mm Fp > 230°F(110°C) Beil. 7(4),2754 FT-NMR 1(1),718A	1g 10g	25.30 131.35
17,976-0 *	FT-IR 1(1),429A SI 71,C,4 R&S 1(1),495E 2-Acetylcyclohexanone, 97% [874-23-7] CH,COC,H,(= O) FW 140.18 bp 111-112°/18mm n\(\tilde{c}\) 1.5090 d 1.078 Fp 175°F(79°C) Beil. 7,559 FT-NMR 1(1),700C FT-IR 1(1),428B SI 69,B,7 Safety 2,36B R&S 1(1),485J	5g 25g	15.70 42.40
	1-Acetyl-1-cyclohexene, 97% [932-66-7] C,H ₅ COCH, FW 124-16 b) 201-202		22.80 73.90
	2-Acetyl-1,3-cyclopentanedione, 99% [3859-39-0] CH ₃ COC ₅ H ₅ (= O) ₂ FW 140.14		16.05
17,977·9 *	2-Acetylcyclopentanone, 98% [1670-46-8] CH ₂ COC ₂ H ₂ (= 0) FW 125.16	100g	11.25 29.45 93.75
36,334-0	N-Acetylcysteamine, 95% [1190-73-4] [N-(2-mercaptoethyl)acetalmide]	1g 5g	24.10 80.30
13,806 -	Chem. Soc., Perkin Trans. 1 1988, 2345. (2) 18traneuri Lett. 166, etc., 166, etc., 166, etc., 178, etc., 186,	10g 50g	11.90 32.60
86,082- 37 791-	(1) J. Chem. Soc., Perkin Irans. 1 1991, 324 l. (2) Telianetto-to-to-to-to-to-to-to-to-to-to-to-to-		28.90 78.65 18.35 63.20
cı 🛴	FT.NMR 1(3),225C SI 391,D,8 R&S 1(2),2477L IRRITANT NH - C - CH ₃ C - CH ₃ HOCH ₂ NH - C - CH ₃	25g	°0

24,707-3

21,467-1

86,082-4

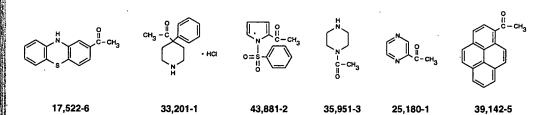
37,791-0

	Acetyldich	\$ 10.00
43,950-9 3-Acetyl-2,5-dichlorothiophene, 98% [36157-40-1] (2,5-dichloro-3-thienyl	1g 5g	10.00 39.00
5-Acetyl-10,11-dihydro-5H-dibenzib,/jazepine, see 30,525 9, 5-Acetyl-10,11-dihydro-5H-dibenzib,/jazepine, see 30,525 9,		22.00
36,921-7 4-Acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-di-2-glass 1(2),2063B FW 216.24 mp 59-61° FT-NMR 1(2),1444B SI 327, A,8 R&S 1(2),2063B FW 216.24 mp 59-61° FT-NMR 1(2),1444B SI 327, A,8 R&S 1(2),2063B 33,227-5 3-O-Acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-di-2-glass 1(2),2063B 33,227-5 3-O-Acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-di-2-glass 1(2),2063B 33,227-5 3-O-Acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-di-2-glass 1(2),2063B FW 216.24 mp 59-61° FT-NMR 1(2),1444B SI 327, A,8 R&S 1(2),2063B 33,227-5 3-O-Acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-di-2-glass 1(2),2063B 3-O-Acetyl-2,5-di-0-lsopropylidene-α-b-glucofuranose, 98% [16713-80-lsopropylidene-α-b-glucofuranose, 98% [170-13-80-lsopropylidene-α-b-glucofuranose, 98% [170-13-80-lsopropylidene-α-b-glucofuran	7] 10g	16.05
FW 302.33 III) 363G RTECS# LZ4910000 S/ 118,C,7 R&S 1(1),763G RTECS# LZ4910000 S/ 118,C,7 R&S 1(1),763G RTECS# LZ4910000	5g	17.20 57.25
cis-3-Acetyl-2,2-dimetriyio) of acid page 1186 acid page 1186 page 122033-07-6] FW 138.17 bp 83°/11mm nB 1.	Ja	18.35 60.70 14.70
0 1,060 FP 100 98% [10599-70-9] FW 15T 18 1/3\ 1457D S/ 364,B	1.4850 5g 3,6 25g	48.85
0 1.030 FP 117 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	21.2//	19.70
IRRITANT [2386-33-6] FW 181.1	19 50g	12.50 41.35 9.75
mp 205-207 Bolimethylthiazole, 96% [38205-60-6] FW 135.25 S. Acetyl-2, 4-dimethylthiazole, 96% [38205-60-6] FW 135.25 S. Acetyl-2, 4-dimethylthiazole, 96% [38205-60-6] FW 135.25 S. Acetyl-2, 5-Acetyl-2, 5-Acety	37D 39	30.05
R&S 1(2),2389E IRRITANT R&S 1(2),2389E IRRITANT R&S 1(2),2389E IRRITANT R&S 1(2),2389E IRRITANT R&S 1(2),2589E IRRITANT R&S 1(15mm	31.30 102.70
10 10 10 10 10 10 10 10 10 10 10 10 10 1		21.80 70.20
Si 368,D,2 Safety 2,38A Nas (G) (22293-38-5] CH,CON(C,H,JNTO,N-2) (29,622-8 N-Acetyl-1,2-diphenylhydrazine, 99% [22293-38-5] CH,CON(C,H,JNTO,N-2) (29,622-8 N-Acetyl-1,2-diphenylhydrazine, 99% [22293-38-5] CH,CON(C,H,JNTO,N-2) (22,034-38-5) CH,CON(C,H,JNTO,N-2) (22,034-38-6) CH,CON(C,H,JNTO,N-2)	rck 17L	97.10
For calibrating analytical instruments.		97.10
Control valve Z16,699-5 or regulator Z16,700-2 is recommended. 32,144-3 Acetylene, 1% in nitrogen [74-86-2] HC=CH	17L	66.85
32,099-4 Acetylene, 1000 ppm in nitrogen [74-6-7-6-7-6-7-6-7-6-7-6-7-6-7-6-7-6-7-6		66.85
32,074-9 Acetylene, 100 ppm in nitrogen [/4-60-2] Tro- For calibrating analytical instruments. Control valve Z16,699-5 or regulator Z16,700-2 is recommended. (Packaged in lightweight no-return cylinder)	:	·

FOR LABORATORY SUPPLIES SEE THE TECHWARE SECTION



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17,522-6	2-Acetylphenothiazine, 95% [6631-94-3] FW 241.31 mp 180-185° FT-NMR 1(2),871A FT-IR 1(2),49B SI 261,D,5 R&S 1(2),1657E	25g 100g	23.35 64.70
39,381-9	4-Acetylphenoxyacetic acid, 99% [1878-81-5] CH ₃ COC ₆ H ₄ OCH ₂ CO ₇ H FW 194.19 mp 175-177° Bell. 8,4,347 SI 280,B,5 IRRITANT	5g 25g	20.95 69.85
19,065-9 ★	α-Acetylphenylacetonitrile, 98% [4468-48-8] (2-phenylacetoacetonitrile)	5g 25g	9.50 31.55
85,745-9 ★	N-Acetyl-L-phenylalanine, 99% [2018-61-3] C₀H₀CH₂CH(NHCOCH₃)CO₂H FW 207.23 mp 171-173° [α]²² + 40.0° (c = 1, CH₃OH) FT-IR 1(2),260B SI 297,A,2 R&S 1(2),1851F	1g 5g	8.45 23.45
85,674-6	N-Acetyl-L-phenylalanyl-3,5-dilodo-L-tyrosine [3786-08-1]	50mg 250mg	11.65 34.45
44,260-7	N-(4-Acetylphenyl)-2-chloroacetamide, 99% [38283-38-4] CICH ₂ CONHC ₆ H ₄ COCH ₃ FW 211.65 mp 153-155° <i>IRRITANT</i>	5g 25g	22.50 75.00
24,271-3 *	1-Acetyl-2-phenylhydrazine, 98% [114-83-0] CH ₃ CONHNHC ₆ H ₆ FW 150.18mp 128-131° Beil. 15,241 FT-NMR 1(2),1413C FT-IR 1(2),356A SI 324;D,5 Safety 2,46C R&S 1(2),2039M RTECS# AJ2900000 TOXIC	25g 100g	16.40 43.80
43,994-0 Œ₩⊅	3-Acetylphenyl isocyanate, 99% [23188-64-9] CH,COC,H,NCO FW 161.16mp 33-34° bp 155°/4mm ng 1.5630 d 1.174 Fp >230°F(110°C) S/ 340,D,9 LACHRYMATOR MOISTURE-SENSITIVE	1g 10g	8.00 45.00
33,201-1	4-Acetyl-4-phenylpiperidine hydrochloride, 98% [10315-03-4] FW 239.75mp 232-234° Beil. 21(3),3702 FT-NMR 1(2),799B SI 253,C,1 R&S 1(2),1619F	1g 5g	26.50 88.20
43,881-2	2-Acetyl-1-(phenylsulfonyl)pyrrole, 97% [86688-88-2] FW 249.29 mp 95-96°	1g 5g	15.00 50.00
43,882-0	3-Acetyl-1-(phenylsulfonyl)pyrrole, 98% [81453-98-7] FW 249.29 mp 96-99°	. 1g 5g	15.00 50.00
	4-Acetylphenyl triflate, see 42,411-0, 4-Acetylphenyl trifluoromethane- sulfonate page 24		•
42,411-0 Œ	4-Acetylphenyl trifluoromethanesulfonate, 99% [109613-00-5] (4-acetylphenyl triflate) CF ₃ SO ₃ C ₄ H ₄ COCH ₃ FW 268.21 bp 75-76°/0.35mm ng 1.4700 d 1.418 Fp > 230°F(110°C) SI 349,D,6 MOISTURE-SENSITIVE TOXIC	5mL 25mL	13.50 45.00
86,214-2	Acetyl phosphate, lithium potassium salt, 97% [94249-01-1] CH,CO ₂ P(O)(OLi)OK FW 184.06 Safety 2,46D R&S 1(1),1121D HYGROSCOPIC	1g 5g	30.75 102.25
35,951-3	1-Acetylpiperazine, 99% [13889-98-0] FW 128.18 mp 32-34° Fp >230°F(110°C) Beil. 23(3),201 FT-NMR 1(1),1250A SI 138,B,1 R&S 1(1),897A IRRITANT HYGROSCOPIC	5g 25g	30.55 106.15
38,825-4	1-Acetyl-4-piperidone, 94% [32161-06-1] FW 141.17 bp 218° ng 1.5030 d 1.146 Fp > 230°F(110°C) Beil. 21,3,3191 FT-NMR 1(1),1251C FT-IR 1(1),765C S/ 138,D,2 R&S 1(1),897I	1mL 5mL	12.15 40.50
26,947-6	N-Acetylprocainamide, 99 + % [32795-44-1]	250mg 1g	14.15 39.65
20,565-6	N-Acetylprocainamide hydrochloride, 99% [34118-92-8]	1g 5g	37.95 109.70
A2,080-4	3-Acetyl-1-propanol, 95% [1071-73-4] (5-hydroxy-2-pentanone)	25g 100g	26.10 71.70
25,180-1 *	Acetylpyrazine, 97% [22047-25-2] FW 122.13 mp 76-78° Beil. 24(3),243	250mg 1g 5g	9.45 26.25 103.95
39,142-5	1-Acetylpyrene, 97% [3264-21-9] FW 244.30 mp 86-502 Bell. 7,3,2726 SI 255,D,8	1g 10g	9.55 53.15



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		Acetylpy	
	2-Acetylpyridine, 99 + % [1122-62-9] FW 121.14 bp 188-189° nB 1.5210 d 1.080	25g 100g	\$ 21.65 77.70
	EN 164 PI/3 C) DOIL - 12 C	10g	12.60
A2,120·7	R&S 1(2),25330 RTECS# OB5310000 IRRITAN1 3-Acetylpyridine, 99% [350-03-8] FW 121.14 bp 220° nB 1.5340 d 1.102	100g	71.10 219.30
A2,140-1	######################################	10g 100g	11.30 40.60
	Fp > 230°F(110°C) Bell. 21,279 F1-NMH (3),3000 IRRITANT Safety 2,48B R&S 1(2),2535L RTECS# OB5426000 IRRITANT Safety 2,48B R&S 1(2),2535L RTECS# OB5426000 IRRITANT 2-Acetylpyrrole, 99% [1072-83-9] (methyl 2-pyrrolyl ketone) FW 109.13 mp > 85° Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(2),567D SI 360,B,4 R&S 1(2),2303N Bell. 21,271 FT-NMR 1(3),5B FT-IR 1(5g 25g	20.15 64.50
	Bell. 21,271 FI-NMR (3,38 PT 17 H 10) RTECS# OB5970000 IRRITANT 5-Acetylsalicylamide, 98% [40187-51-7] (5-acetyl-2-hydroxybenzamide) 5-Acetylsalicylamide, 98% [40187-51-7] (5-acetyl-2-hydroxybenzamide) 5-Acetylsalicylamide, 98% [40187-51-7] (5-acetyl-2-hydroxybenzamide) 5-Acetylsalicylamide, 98% [40187-51-7] (5-acetyl-2-hydroxybenzamide)	25g 100g	25.65 84.30
	SI 323, B.4 R&S 1(2), 2025J RTECS# CU8702280 IRRITANT	1g	7.70 23.25
*	Acetylsalicylic acid, 99 + % [50-76-2] (2006 Bell. 10,67 Merck Index 11,873 2-(CH,CO,)C,H,CO,H FW 180.16 mp 138-140° Bell. 10,67 Merck Index 11,873 2-(CH,CO,H,CO,H FW 180.16 mp 138-140° Bell. 10,167 Merck Index 11,873 2-(CH,CO,H,CO,H,CO)		5.10
13,292-6 *	ACETylsalicylic acid, 99% [50-78-2] (2-acetoxybenzoic acid) 2-(CH ₃ CO ₃ C ₄ H ₄ CO ₃ H ₄ CO ₄ H ₄ CO ₄ H ₄ CO ₄ H ₄ CO ₅ H ₅ CO ₅ H ₄ CO	-	18.65 57.35
16,519-0	Acetylsalicyloyl chloride, 95% [5538-51-2] 2-(CH,CO ₂)C ₆ H ₆ COCI FW 198.61	. 5g 25g	16.60 66.85
	Used for adviation of cyclobation at A annual 5 hydroxytryntamine page 21	100g	17.40
29,272-	N-Acetylserotonin, see 85,548-0, N-Acetyl-3-11940-07417948-08-1940-07417948-0741949-08-1949-0	500g	57.70
•	HYGROSCOPIC	100g 500g	14.30 32.10
	Chloride) 4-(Ch. Cotti, Ja. 102, 1622C FT-IR 1(2), 522C ST 350, 0,4	•	45.70 35.05
A2,220	Safety 2,50B H&S 1(2),2231 (CH,CO) ₂ S FW 118.15 bp 119-120° nB 1.4790 d 1.117		114.55
	R&S 1(1),829A FLAMMABLE LINE 00 00 TW 100 23 mp 55-57° bp 135-142°/1mm	5g 25g	34.60 118.50
15,037 41,319	2-Acetyl-1-tetralone, 98% [17276-08-9] FW 188.25 III. 166.25 III.	50mL 250mL	13.80 46.65
28.84°	CORROSIVE 2-Acetylthlazole, 99% [24295-03-2] FW 127.17 bp 89-91°/12mm ng 1.5480 d 1.227 Fp 173°-[78°C) Bell. 27(3),2617 FT-IR 1(3),1494D S/ 379,A,4 Safety 2,50D	250mg 1g 5g	10.35 28.65 91.30
,	R&S 1(2),2389D STENCH Useful for preparation of triazolothiazoles, chiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and in aldol useful for preparation of triazolothiazoles, thiral alcohols, and	·	9.50
04.05	1991, 2, 243. (3) J. Org. Cristil. 1801, 65, 65	1g 5g	32.00
34,95 ••••• 85,53	3.2 S-Acetylthiocholine bromide, 98% [25025-59-6] [(2-mercaptoethyl)trimethyl	1g 5g	20.65 74.00
	mp 217-223° (dec.) FI-IN 1(1),0100 G. 1-259-59 HYGROSCOPIC		
	CH ₃ (CH ₂) ₁₀ CH ₂ CH ₃ CH ₃ (CH ₂) ₁₀ CH ₂ CH ₃	[s]	о - -сн _а
 A0 44	no-2 24,735-9 15,037-1 41,319-4	28,84	i 1-1
A2,10	[-N		

1 15	





■ Acetylthio ■

Acety			\$
A2,230-0 *	S-Acetylthiocholine iodide, 98% [1866-15-5] [(2-mercaptoethyl)trimethyl	1g 5g	9.90 33.00
A2,260-2	2-Acetylthlophene, 98% [88-15-3] (methyl 2-thlenyl ketone) FW 126.18 mp 10-11°	25g 100g	8:50 21.95
19,632-0	bp 214° nB 1.5650 d 1.168 Fp 196°F(31°C) Bell. (1,02) ABC 12,535N RTECS# OB6300000 FT-IR 1(2),596C SI 368,C,1 Safety 2,51C R&S 1(2),2335N RTECS# OB6300000 ST-IR 1(2),5960 SI 368,C,1 Safety 2,51C R&S 1(2),2335N RTECS# OB6300000 FT-IR 1(2),597B mp 59-63° bp 208-210/7/48mm Bell. 17(3),4520 FT-NMR 1(3),54B FT-IR 1(2),597B	1g 10g	11.65 63.40
A2,280·7	S/368,C,2 Safety 2,510 H&S 1(2,2337) 1-Acetyl-3-thiosemicarbazide, 95% [2302-88-7] CH,CONHNHCSNH, FW 133.17 1-Acetyl-3-thiosemicarbazide, 95% [2302-88-7] CH,CONHNHCSNH, FW 133.17	25g	30.95
	2-(Acetylthio)succinic anhydride, see 19,732-7, S-Acetylthiercaptosuccinic	25a	10.90
A2,285-8 ★	1.Acetyl-2-thiourea, 99% [591-08-2] CH,CONHCSNH, FW 118.16 mp 165-169°	100g	29.30
	a a real accompany soo 24 917.7 Vitamin E acetate page 1551		28.85
34,727-2	1-Acetyl-1H-1,2,3-triazolo[4,5-b]pyridine, 97 % [707800-34-0] FW 102-13	1g 5g	95.20
•	1-O-Acetyl-2,3,5-tril-O-benzoyl-β-p-riboturanose, see 15,501-6, β-p-riboturanose,	1g	11.95
36,282-4	Acetyltrimethylsilane, 97% [13411-48-8] (CH ₃),SICOCH ₃ FW 116.24 7/6 1.3 100.811 Fp 48°F(8°C) FT-MRI (3),682A SI 472,D,5 R&S 1(2),2983A SI (2),2983A SI (2),29	5g 25g	44.00 146.90
	Useful as a hindered acetaldehyde equivalent in stereo-commoned and reasons and the state of the	5g	8.40
85,580-4 *	N-Acetyl-pt-tryptophan, 99 + % [87-32-1] FW 246.27 mp 204-206° (dec.) FT-NMR 1(3),147A FT-IR 1(2),674B SI 383,E,4 R&S 1(2),2415K	25g 100g	26.55 81.90
	N-Acetyl-L-tryptophanamide, 98% [2382-79-8] FW 245.28 mp 194-196°	250mg	17.90
85,675-4	$[\alpha]^{22} + 17.5^{\circ} (c = 2, CH_3OH)$ F1-NMH $I(O)$, 140A 77 H $I(O)$		46.25 9.75
85,772-6	N-Acetyl-L-tryptophan ethyl ester, 99% [2382-80-1] FW 274.32 mp 112-114°		30.70 18.80
85,531-6	N-Acetyl-L-tyrosinamide, 99 % [1948-71-6] 4-(HO)C ₆ H ₂ CH ₂ CH ₃		18.95
44,153-8 Φ ★	N-Acetyl-L-tyrosine, 98% [537-55-3] 4-(HO)C ₄ H ₄ CH ₂ CH(NHCOCH ₃)CO ₂ H FW 223.23		83.80 9.40
A2,290-4	N-Acetyl-L-tyrosine ethyl ester monohydrate, 99% [36546-50-6] (A1EC)	1g 10g	37.50
21,100-1	[α]β + 24.1° (c = 1, C ₂ H ₂ OH) F1-IR 1(2),60°B 31-23(3,4) No. (c),10°C (c	. 25g	11.50
•	Dye content ~50% Acid Black 1, see 19,524-3, Naphthol Blue Black page 1058		
	A LLA PLACE A COO 10 828-5 Nigrosin water soluble page 1071		-4.45
21,045-	5 Acid Black 24 [3071-73-6] (C.I. 26370) FW 731.72 xmax 5721111 F1-17 (2),0025	. 50g	21.10
•	Dye content ~50%		
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A2,260-2

34,727-2

85,580-4

85,675-4

21,100-1

NaO₃S - N = N - NH - SO₃Na

21,045-5